

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122623\
 Data File : BM043595.D
 Acq On : 27 Dec 2023 04:58
 Operator : MA/JU
 Sample : PB158081BS
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS081

Quant Time: Dec 27 05:29:09 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM121123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 26 22:58:31 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.971	152	5340	0.400	ng/ul	0.00
4) Naphthalene-d8	10.779	136	15570	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.606	164	6448	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.350	188	12380	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.527	240	7596	0.400	ng/ul	# 0.00
23) Perylene-d12	23.939	264	8562	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.376	96	5252	0.778	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.368	152	7236	0.328	ng/ul	0.00
18) Fluoranthene-d10	19.371	212	9329	0.344	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.410	88	15917	2.352	ng/ul#	81
5) Naphthalene	10.829	128	16215	0.346	ng/ul	99
7) 2-Methylnaphthalene	12.440	142	9393	0.311	ng/ul	99
8) 1-Methylnaphthalene	12.654	142	9072	0.299	ng/ul	100
10) Acenaphthylene	14.328	152	12696	0.358	ng/ul	98
11) Acenaphthene	14.666	153	9252	0.358	ng/ul	98
12) Fluorene	15.656	166	9865	0.337	ng/ul	99
14) Pentachlorophenol	16.999	266	2113	0.770	ng/ul	99
15) Phenanthrene	17.392	178	14720	0.357	ng/ul	98
16) Anthracene	17.485	178	13815	0.368	ng/ul	98
19) Fluoranthene	19.403	202	14464	0.338	ng/ul#	92
20) Pyrene	19.766	202	14899	0.350	ng/ul#	94
21) Benzo(a)anthracene	21.512	228	11873	0.356	ng/ul	99
22) Chrysene	21.565	228	13295	0.378	ng/ul	99
24) Benzo(b)fluoranthene	23.202	252	13189	0.339	ng/ul	91
25) Benzo(k)fluoranthene	23.254	252	14812	0.367	ng/ul#	88
26) Benzo(a)pyrene	23.833	252	12285	0.383	ng/ul#	86
27) Indeno(1,2,3-cd)pyrene	26.455	276	16312	0.447	ng/ul#	90
28) Dibenzo(a,h)anthracene	26.485	278	13182	0.456	ng/ul#	93
29) Benzo(g,h,i)perylene	27.219	276	13913	0.446	ng/ul	95

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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